



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:15 PM GMT

PDB ID : 1JH0  
Title : Photosynthetic Reaction Center Mutant With Glu L 205 Replaced to Leu  
Authors : Camara-Artigas, A.; Magee, C.L.; Williams, J.C.; Allen, J.P.  
Deposited on : 2001-06-27  
Resolution : 3.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

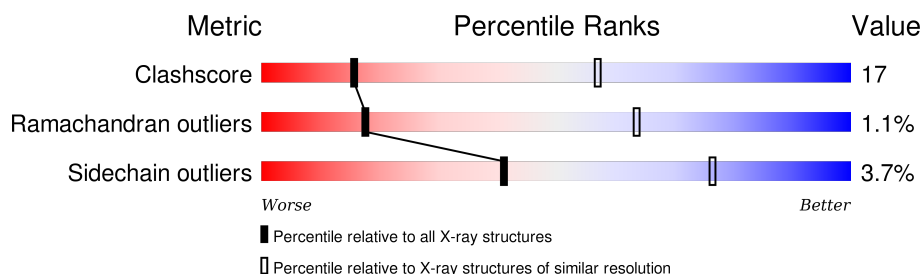
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	281	 64% 33% •
2	M	307	 68% 28% • •
3	H	260	 67% 23% • 8%

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 7025 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Photosynthetic Reaction Center L subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	280	Total	C	N	O	S	0	0	0
			2226	1506	354	358	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	205	LEU	GLU	ENGINEERED	UNP P02954

- Molecule 2 is a protein called Photosynthetic Reaction Center M subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	301	Total	C	N	O	S	0	0	0
			2404	1605	393	396	10			

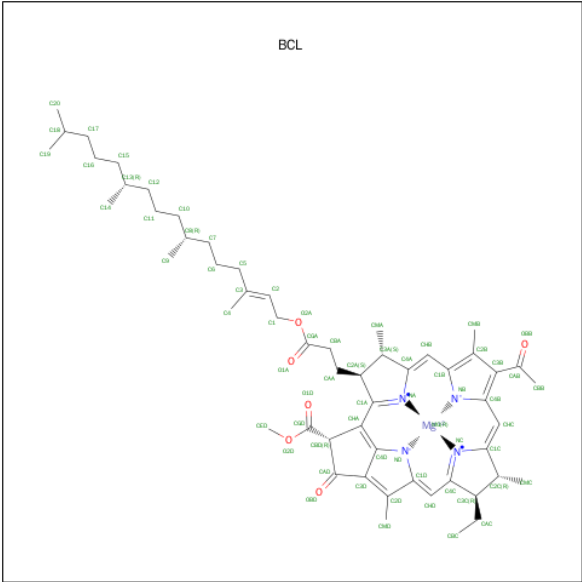
- Molecule 3 is a protein called Photosynthetic Reaction Center H subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	238	Total	C	N	O	S	0	0	0
			1814	1160	311	334	9			

- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

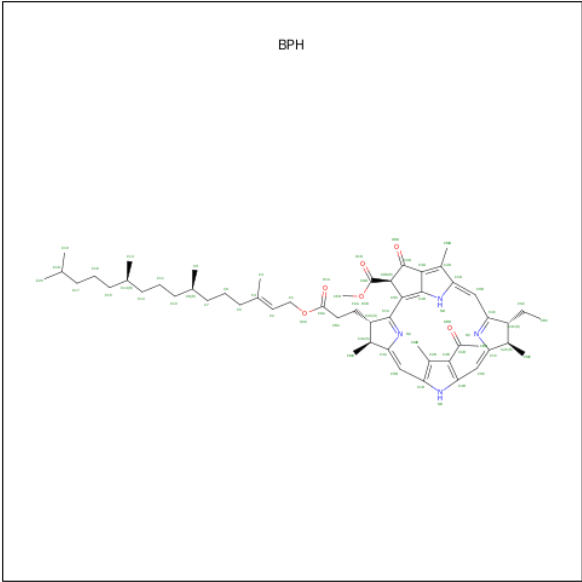
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	M	1	Total	Fe	0	0
			1	1		

- Molecule 5 is BACTERIOCHLOROPHYLL A (three-letter code: BCL) (formula: C<sub>55</sub>H<sub>74</sub>MgN<sub>4</sub>O<sub>6</sub>).



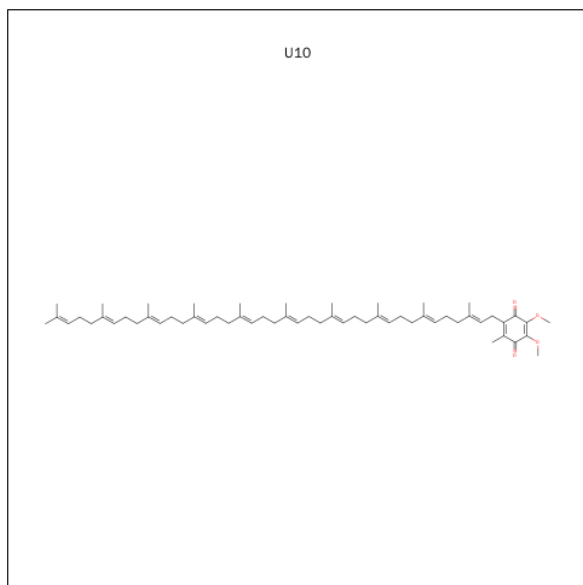
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 6 is BACTERIOPHEOPHYTIN A (three-letter code: BPH) (formula: C<sub>55</sub>H<sub>76</sub>N<sub>4</sub>O<sub>6</sub>).



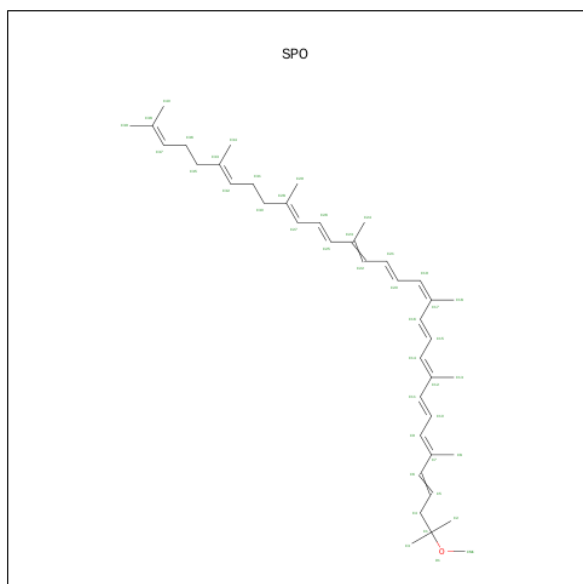
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	M	1	Total	C	N	O	0	0
			65	55	4	6		
6	L	1	Total	C	N	O	0	0
			65	55	4	6		

- Molecule 7 is UBIQUINONE-10 (three-letter code: U10) (formula:  $C_{59}H_{90}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	M	1	Total	C	O	0	0
			48	44	4		

- Molecule 8 is SPHEROIDENE (three-letter code: SPO) (formula:  $C_{41}H_{60}O$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	M	1	Total	C	O	0	0
			42	41	1		

- Molecule 9 is water.

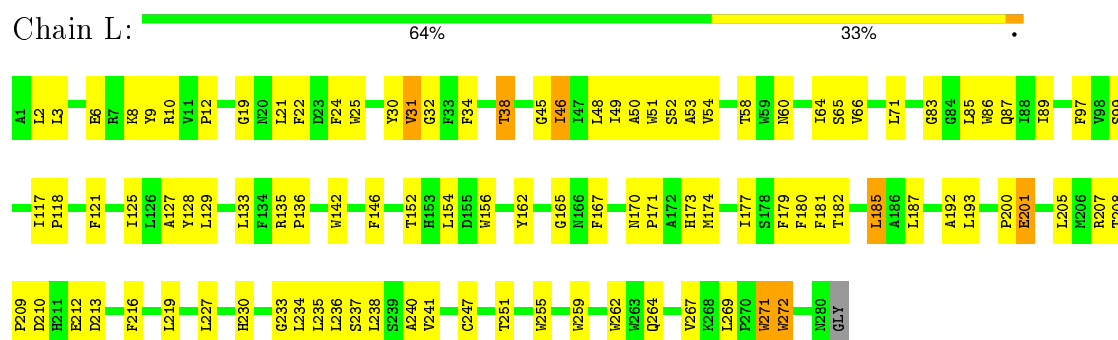
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	H	28	Total	O	0	0
			28	28		
9	L	23	Total	O	0	0
			23	23		
9	M	45	Total	O	0	0
			45	45		

### 3 Residue-property plots

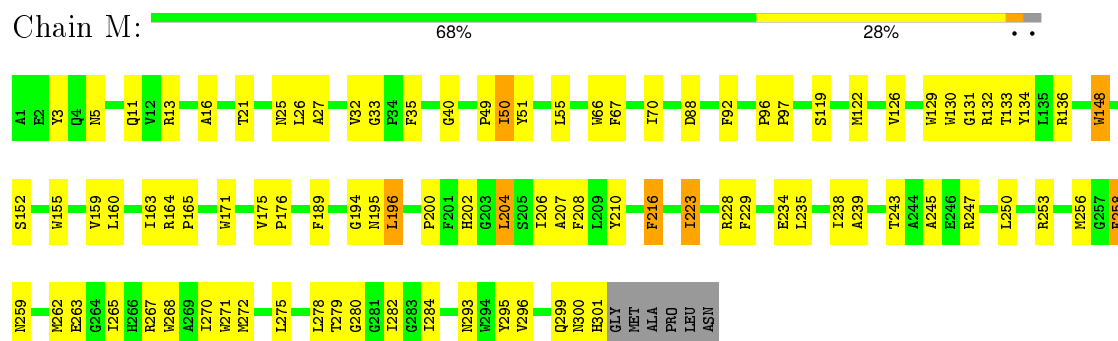
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

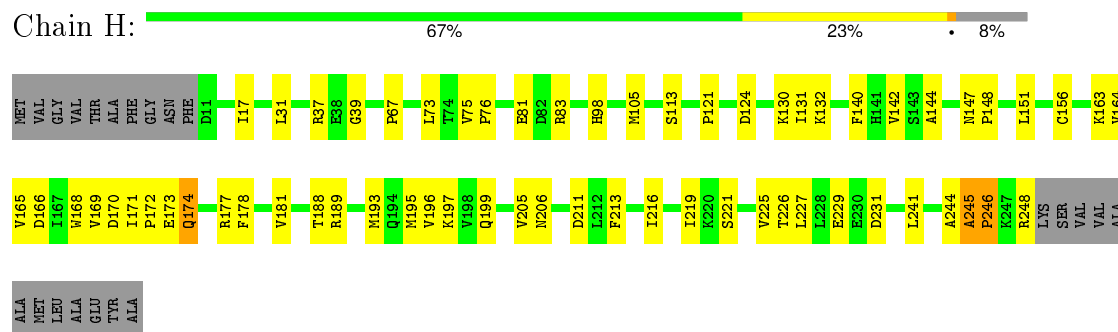
#### • Molecule 1: Photosynthetic Reaction Center L subunit



#### • Molecule 2: Photosynthetic Reaction Center M subunit



#### • Molecule 3: Photosynthetic Reaction Center H subunit



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	141.80 Å   141.80 Å   187.40 Å 90.00°   90.00°   120.00°	Depositor
Resolution (Å)	29.90 – 3.50	Depositor
% Data completeness (in resolution range)	99.5 (29.90-3.50)	Depositor
$R_{merge}$	0.23	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.225 , 0.269	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7025	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BCL, BPH, U10, FE, SPO

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	L	0.44	0/2314	0.64	0/3169
2	M	0.45	0/2496	0.62	0/3408
3	H	0.44	0/1862	0.69	0/2534
All	All	0.45	0/6672	0.65	0/9111

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	2226	0	2189	91	0
2	M	2404	0	2318	91	0
3	H	1814	0	1818	56	0
4	M	1	0	0	0	0
5	L	132	0	148	14	0
5	M	132	0	148	17	0
6	L	65	0	76	8	0
6	M	65	0	76	9	0
7	M	48	0	63	3	0
8	M	42	0	60	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	H	28	0	0	0	0
9	L	23	0	0	3	0
9	M	45	0	0	1	0
All	All	7025	0	6896	236	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (236) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:48:LEU:HB2	1:L:89:ILE:HD11	1.37	1.04
3:H:156:CYS:SG	3:H:248:ARG:HB2	2.13	0.88
1:L:97:PHE:HB3	1:L:125:ILE:HD12	1.59	0.83
1:L:58:THR:HG23	1:L:64:ILE:HD13	1.60	0.82
1:L:48:LEU:HB2	1:L:89:ILE:CD1	2.12	0.79
3:H:248:ARG:HB2	3:H:248:ARG:NH1	1.98	0.79
1:L:48:LEU:CB	1:L:89:ILE:HD11	2.12	0.79
1:L:117:ILE:HD12	1:L:117:ILE:H	1.47	0.78
3:H:196:VAL:HG12	3:H:205:VAL:HG22	1.66	0.78
1:L:241:VAL:HG21	6:L:855:BPH:HAC2	1.70	0.73
1:L:219:LEU:HD12	2:M:132:ARG:NH1	2.04	0.73
2:M:238:ILE:HD13	2:M:262:MET:HB3	1.73	0.70
3:H:130:LYS:HE3	3:H:170:ASP:OD2	1.92	0.70
6:L:855:BPH:HBB2	2:M:210:TYR:HB3	1.73	0.70
3:H:148:PRO:HA	3:H:151:LEU:HD12	1.72	0.70
3:H:132:LYS:HD2	3:H:171:ILE:HD11	1.73	0.69
1:L:173:HIS:CE1	1:L:177:ILE:HD11	2.27	0.69
1:L:192:ALA:HB2	2:M:270:ILE:HD11	1.75	0.68
1:L:97:PHE:HB3	1:L:125:ILE:CD1	2.24	0.68
1:L:205:LEU:HD13	3:H:67:PRO:HA	1.76	0.68
1:L:219:LEU:HD11	2:M:133:THR:HG22	1.74	0.68
1:L:174:MET:HB3	5:M:850:BCL:O1D	1.94	0.67
1:L:38:THR:HG22	1:L:99:SER:CB	2.24	0.67
3:H:156:CYS:SG	3:H:248:ARG:CB	2.82	0.67
2:M:245:ALA:HB2	2:M:262:MET:HE2	1.75	0.67
2:M:196:LEU:HD23	2:M:202:HIS:CD2	2.30	0.66
3:H:142:VAL:HG21	3:H:147:ASN:ND2	2.10	0.66
1:L:51:TRP:O	1:L:54:VAL:HG22	1.95	0.66
1:L:38:THR:HG22	1:L:99:SER:HB2	1.75	0.66
3:H:248:ARG:HH11	3:H:248:ARG:HB2	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:135:ARG:HB3	1:L:136:PRO:HD3	1.78	0.65
1:L:264:GLN:HA	1:L:267:VAL:HG12	1.79	0.64
2:M:253:ARG:HB2	2:M:259:ASN:HD22	1.61	0.64
2:M:200:PRO:HB2	3:H:17:ILE:HD12	1.79	0.63
1:L:271:TRP:CD1	1:L:271:TRP:N	2.65	0.62
3:H:156:CYS:SG	3:H:248:ARG:HA	2.39	0.62
1:L:60:ASN:O	1:L:64:ILE:HG12	2.00	0.62
2:M:228:ARG:NE	3:H:195:MET:HE3	2.13	0.62
1:L:121:PHE:CZ	1:L:125:ILE:HD11	2.34	0.62
1:L:181:PHE:HB3	6:M:854:BPH:CBB	2.28	0.62
1:L:6:GLU:HG3	2:M:250:LEU:HD21	1.80	0.62
1:L:45:GLY:HA3	6:L:855:BPH:H9C1	1.82	0.62
2:M:16:ALA:HB1	2:M:32:VAL:HG21	1.82	0.61
2:M:119:SER:HB3	8:M:859:SPO:H311	1.83	0.61
3:H:165:VAL:O	3:H:166:ASP:HB2	2.00	0.60
3:H:131:ILE:HD13	3:H:170:ASP:HA	1.83	0.60
1:L:182:THR:HG22	1:L:236:LEU:HD13	1.83	0.60
1:L:34:PHE:O	1:L:38:THR:HG23	2.02	0.60
2:M:159:VAL:HA	2:M:163:ILE:HB	1.84	0.59
1:L:192:ALA:N	2:M:270:ILE:HD13	2.17	0.59
5:M:850:BCL:H11	6:M:854:BPH:HBB2	1.84	0.58
3:H:131:ILE:CD1	3:H:177:ARG:HD2	2.33	0.58
3:H:156:CYS:SG	3:H:248:ARG:CA	2.91	0.58
5:M:850:BCL:HMB2	6:M:854:BPH:HMB3	1.85	0.58
2:M:275:LEU:HD23	2:M:278:LEU:HD23	1.85	0.58
1:L:49:ILE:HD13	1:L:66:VAL:HG21	1.86	0.58
2:M:202:HIS:O	2:M:206:ILE:HD13	2.04	0.57
1:L:192:ALA:HB2	2:M:270:ILE:CD1	2.35	0.57
3:H:130:LYS:HZ1	3:H:172:PRO:HG2	1.70	0.57
3:H:81:GLU:O	3:H:83:ARG:HG2	2.04	0.57
1:L:269:LEU:HB2	1:L:272:TRP:NE1	2.20	0.57
1:L:208:THR:HB	1:L:209:PRO:HD2	1.85	0.56
5:M:850:BCL:CBB	5:M:850:BCL:HHC	2.36	0.56
1:L:52:SER:HB2	1:L:85:LEU:HD13	1.87	0.56
5:L:853:BCL:HMD2	5:M:852:BCL:HBB3	1.87	0.55
2:M:50:ILE:HD13	2:M:51:TYR:C	2.27	0.55
2:M:267:ARG:HG2	2:M:267:ARG:HH11	1.71	0.55
2:M:21:THR:HG23	2:M:26:LEU:HD21	1.88	0.55
1:L:58:THR:HG23	1:L:64:ILE:CD1	2.35	0.54
2:M:267:ARG:O	2:M:270:ILE:HG22	2.07	0.54
2:M:11:GLN:HB2	3:H:144:ALA:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:202:HIS:CE1	2:M:206:ILE:HD11	2.41	0.54
2:M:280:GLY:HA2	5:M:852:BCL:HED2	1.90	0.54
5:M:850:BCL:HBB3	5:M:850:BCL:HHC	1.90	0.54
1:L:255:TRP:CZ2	1:L:262:TRP:HB2	2.43	0.54
3:H:168:TRP:HB2	3:H:178:PHE:HB2	1.90	0.53
2:M:132:ARG:O	2:M:136:ARG:HG2	2.09	0.53
3:H:131:ILE:HD12	3:H:177:ARG:HD2	1.91	0.53
5:L:853:BCL:HAA2	5:L:853:BCL:HBD	1.91	0.53
2:M:267:ARG:HG2	2:M:267:ARG:NH1	2.22	0.53
5:L:851:BCL:H122	6:L:855:BPH:H3A	1.91	0.53
2:M:25:ASN:OD1	2:M:27:ALA:HB3	2.10	0.52
2:M:160:LEU:HD23	2:M:284:ILE:HG21	1.91	0.52
1:L:83:GLY:O	1:L:87:GLN:HG3	2.09	0.52
2:M:164:ARG:HB3	2:M:165:PRO:HD3	1.91	0.52
1:L:117:ILE:HB	1:L:118:PRO:CD	2.40	0.52
3:H:226:THR:OG1	3:H:229:GLU:HG3	2.09	0.52
1:L:181:PHE:CD2	6:M:854:BPH:HBB1	2.45	0.52
2:M:129:TRP:O	2:M:133:THR:HG23	2.10	0.52
2:M:32:VAL:HG12	2:M:33:GLY:O	2.09	0.52
1:L:117:ILE:HB	1:L:118:PRO:HD3	1.93	0.51
3:H:248:ARG:CB	3:H:248:ARG:NH1	2.73	0.51
1:L:49:ILE:CD1	1:L:66:VAL:HG21	2.41	0.51
2:M:253:ARG:HB2	2:M:259:ASN:ND2	2.26	0.51
1:L:87:GLN:NE2	1:L:142:TRP:CD1	2.80	0.50
3:H:213:PHE:O	3:H:216:ILE:HG13	2.12	0.50
1:L:177:ILE:HG12	5:L:851:BCL:HMB3	1.94	0.50
5:L:851:BCL:HBB3	5:M:850:BCL:HMD2	1.94	0.50
2:M:234:GLU:O	2:M:238:ILE:HG12	2.12	0.50
1:L:50:ALA:O	1:L:53:ALA:HB3	2.12	0.50
3:H:241:LEU:O	3:H:248:ARG:NH2	2.45	0.50
3:H:244:ALA:O	3:H:246:PRO:HD2	2.11	0.50
3:H:219:ILE:HD12	3:H:221:SER:O	2.12	0.50
1:L:181:PHE:HB3	6:M:854:BPH:HBB2	1.93	0.49
5:L:851:BCL:CAA	5:L:853:BCL:HBC1	2.42	0.49
2:M:155:TRP:HZ2	2:M:282:ILE:HD13	1.78	0.49
3:H:193:MET:O	3:H:196:VAL:HG22	2.13	0.49
1:L:12:PRO:HD3	3:H:98:HIS:O	2.11	0.49
1:L:6:GLU:HG3	2:M:250:LEU:CD2	2.42	0.49
1:L:179:PHE:HB3	1:L:240:ALA:HB2	1.93	0.49
2:M:50:ILE:HD13	2:M:51:TYR:N	2.28	0.49
5:L:851:BCL:HAA2	5:L:853:BCL:HBC1	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:206:ILE:HG13	5:M:852:BCL:HMB3	1.95	0.48
3:H:151:LEU:O	3:H:164:VAL:HG23	2.14	0.48
1:L:46:ILE:N	1:L:46:ILE:HD13	2.27	0.48
2:M:256:MET:CE	2:M:258:PHE:CE2	2.96	0.48
1:L:31:VAL:HG12	1:L:32:GLY:N	2.29	0.48
1:L:34:PHE:HB2	9:L:939:HOH:O	2.13	0.48
1:L:2:LEU:HD21	1:L:10:ARG:CZ	2.43	0.48
2:M:245:ALA:HB2	2:M:262:MET:CE	2.40	0.48
9:L:897:HOH:O	2:M:49:PRO:HG2	2.14	0.48
3:H:75:VAL:HA	3:H:76:PRO:C	2.35	0.48
1:L:52:SER:HB2	1:L:85:LEU:CD1	2.44	0.47
2:M:96:PRO:HB2	2:M:97:PRO:HD2	1.95	0.47
1:L:219:LEU:HA	2:M:132:ARG:HH12	1.79	0.47
3:H:245:ALA:HB3	3:H:246:PRO:HD3	1.96	0.47
2:M:97:PRO:CG	2:M:171:TRP:HB2	2.44	0.47
2:M:239:ALA:O	3:H:73:LEU:HD22	2.14	0.47
1:L:6:GLU:OE2	1:L:10:ARG:NH2	2.46	0.47
2:M:206:ILE:HG23	5:M:852:BCL:HMB3	1.96	0.47
5:M:852:BCL:H201	6:M:854:BPH:H8	1.96	0.47
1:L:25:TRP:CD1	1:L:30:TYR:HA	2.50	0.47
2:M:13:ARG:O	3:H:140:PHE:HA	2.14	0.47
2:M:3:TYR:CZ	2:M:5:ASN:HA	2.50	0.47
2:M:228:ARG:CD	3:H:195:MET:HE3	2.44	0.47
2:M:97:PRO:HG2	2:M:171:TRP:HB2	1.95	0.47
3:H:130:LYS:NZ	3:H:172:PRO:HG2	2.29	0.46
3:H:131:ILE:HD11	3:H:177:ARG:HD2	1.97	0.46
3:H:131:ILE:HA	3:H:169:VAL:O	2.15	0.46
3:H:226:THR:O	3:H:227:LEU:C	2.53	0.46
1:L:171:PRO:HD2	1:L:259:TRP:CZ3	2.51	0.46
1:L:38:THR:HG22	1:L:99:SER:HB3	1.95	0.46
1:L:180:PHE:CD2	1:L:240:ALA:HB1	2.51	0.46
2:M:194:GLY:O	2:M:195:ASN:HB3	2.15	0.46
1:L:117:ILE:H	1:L:117:ILE:CD1	2.24	0.46
3:H:248:ARG:CB	3:H:248:ARG:CZ	2.94	0.46
5:M:850:BCL:HBD	5:M:850:BCL:HAA2	1.98	0.46
1:L:127:ALA:CB	5:L:851:BCL:H43	2.47	0.45
2:M:235:LEU:O	2:M:238:ILE:HB	2.16	0.45
2:M:275:LEU:CD2	2:M:278:LEU:HD23	2.46	0.45
3:H:163:LYS:O	3:H:181:VAL:HG13	2.15	0.45
2:M:66:TRP:CD1	2:M:122:MET:HB2	2.52	0.45
2:M:268:TRP:CE3	3:H:31:LEU:HD13	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:37:ARG:HG2	3:H:37:ARG:HH11	1.81	0.45
2:M:271:TRP:O	2:M:272:MET:C	2.55	0.45
2:M:148:TRP:HA	2:M:148:TRP:CE3	2.51	0.45
2:M:270:ILE:HG23	2:M:271:TRP:N	2.32	0.45
5:M:850:BCL:H2	6:M:854:BPH:HMB2	1.98	0.45
2:M:284:ILE:HG12	5:M:852:BCL:HED3	1.98	0.45
3:H:173:GLU:O	3:H:174:GLN:C	2.54	0.45
1:L:127:ALA:HB1	5:L:851:BCL:H43	1.98	0.45
3:H:156:CYS:HB3	3:H:206:ASN:O	2.17	0.45
1:L:219:LEU:HD12	2:M:132:ARG:HH11	1.80	0.45
1:L:182:THR:OG1	5:M:850:BCL:H42	2.16	0.44
2:M:66:TRP:O	2:M:70:ILE:HG12	2.17	0.44
3:H:248:ARG:O	3:H:248:ARG:HG2	2.15	0.44
5:L:851:BCL:HAA2	5:L:851:BCL:HBD	1.99	0.44
2:M:88:ASP:HB2	2:M:92:PHE:CZ	2.53	0.44
2:M:134:TYR:CD1	2:M:134:TYR:C	2.91	0.44
1:L:170:ASN:O	1:L:174:MET:HG3	2.18	0.44
1:L:128:TYR:HB2	5:L:851:BCL:H61	2.00	0.44
2:M:238:ILE:HD12	2:M:263:GLU:HB2	1.98	0.44
1:L:66:VAL:HG12	1:L:86:TRP:HB2	2.00	0.44
1:L:269:LEU:HD12	1:L:272:TRP:CZ2	2.53	0.44
1:L:162:TYR:HA	1:L:165:GLY:O	2.17	0.44
6:L:855:BPH:HHB	6:L:855:BPH:HBB3	2.00	0.44
1:L:200:PRO:O	1:L:201:GLU:O	2.36	0.44
2:M:293:ASN:OD1	2:M:295:TYR:HB3	2.18	0.44
3:H:189:ARG:HG2	3:H:189:ARG:HH11	1.83	0.43
3:H:197:LYS:HE3	3:H:199:GLN:NE2	2.32	0.43
6:M:854:BPH:HHB	6:M:854:BPH:HMB1	1.78	0.43
2:M:296:VAL:O	2:M:299:GLN:HB2	2.19	0.43
2:M:11:GLN:OE1	2:M:40:GLY:HA3	2.18	0.43
2:M:164:ARG:CZ	2:M:189:PHE:HE2	2.32	0.43
2:M:13:ARG:HD3	2:M:35:PHE:CD2	2.54	0.43
1:L:22:PHE:HA	1:L:24:PHE:CE2	2.53	0.43
1:L:3:LEU:HD22	3:H:39:GLY:HA3	2.00	0.43
2:M:130:TRP:O	2:M:131:GLY:C	2.57	0.42
2:M:275:LEU:HA	2:M:278:LEU:HB3	2.02	0.42
3:H:181:VAL:O	3:H:188:THR:HA	2.19	0.42
2:M:148:TRP:HE3	2:M:148:TRP:HA	1.84	0.42
1:L:146:PHE:HB3	1:L:156:TRP:CD2	2.54	0.42
3:H:124:ASP:C	3:H:124:ASP:OD1	2.57	0.42
1:L:8:LYS:HE2	1:L:9:TYR:CE2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:265:ILE:HG21	7:M:857:U10:C3M	2.48	0.42
7:M:857:U10:H201	7:M:857:U10:H222	1.78	0.42
2:M:55:LEU:HD12	2:M:55:LEU:HA	1.86	0.42
1:L:238:LEU:HD23	6:L:855:BPH:CBC	2.50	0.42
1:L:219:LEU:O	2:M:132:ARG:NH1	2.46	0.42
2:M:152:SER:O	2:M:155:TRP:HB3	2.20	0.42
1:L:193:LEU:HD23	9:L:941:HOH:O	2.19	0.42
1:L:129:LEU:O	1:L:133:LEU:HB3	2.20	0.42
5:L:853:BCL:O1A	5:L:853:BCL:H43	2.19	0.42
1:L:65:SER:CB	1:L:152:THR:HG21	2.50	0.42
1:L:200:PRO:O	1:L:201:GLU:C	2.58	0.42
2:M:175:VAL:HG13	2:M:176:PRO:HD2	2.02	0.42
1:L:230:HIS:CD2	2:M:223:ILE:HG21	2.55	0.42
1:L:251:THR:HG1	1:L:259:TRP:HZ2	1.68	0.41
2:M:204:LEU:O	2:M:207:ALA:HB3	2.19	0.41
1:L:227:LEU:O	1:L:227:LEU:HG	2.20	0.41
2:M:148:TRP:CD1	9:M:901:HOH:O	2.57	0.41
1:L:212:GLU:O	1:L:213:ASP:C	2.58	0.41
1:L:208:THR:C	1:L:210:ASP:N	2.72	0.41
1:L:233:GLY:HA3	2:M:216:PHE:CE1	2.55	0.41
7:M:857:U10:H271	7:M:857:U10:H251	1.81	0.41
1:L:117:ILE:HD12	1:L:117:ILE:N	2.24	0.41
1:L:185:LEU:C	1:L:185:LEU:HD23	2.41	0.41
1:L:187:LEU:HB2	2:M:216:PHE:CD2	2.55	0.41
2:M:247:ARG:NH2	3:H:113:SER:O	2.52	0.41
5:L:851:BCL:H2C	5:L:851:BCL:HBC2	1.95	0.41
6:L:855:BPH:CBB	2:M:210:TYR:HB3	2.47	0.41
5:M:850:BCL:CHC	5:M:850:BCL:HBB3	2.49	0.41
2:M:253:ARG:HH11	2:M:253:ARG:HG3	1.85	0.41
2:M:300:ASN:O	2:M:301:HIS:HB2	2.21	0.41
1:L:45:GLY:HA3	6:L:855:BPH:C9	2.50	0.41
2:M:253:ARG:HG3	2:M:253:ARG:NH1	2.36	0.41
3:H:216:ILE:HG13	3:H:216:ILE:H	1.57	0.41
2:M:208:PHE:HE2	2:M:279:THR:HG21	1.86	0.41
2:M:228:ARG:HG3	2:M:229:PHE:CE2	2.56	0.40
5:L:851:BCL:CBB	5:M:850:BCL:HMD2	2.51	0.40
3:H:142:VAL:CG2	3:H:147:ASN:ND2	2.82	0.40
1:L:269:LEU:HD12	1:L:272:TRP:HZ2	1.85	0.40
1:L:234:LEU:O	1:L:238:LEU:HG	2.22	0.40
2:M:67:PHE:CD1	6:M:854:BPH:H9C1	2.57	0.40
3:H:130:LYS:NZ	3:H:172:PRO:CG	2.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:243:THR:O	2:M:247:ARG:HG3	2.22	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	L	278/281 (99%)	246 (88%)	27 (10%)	5 (2%)	11	53
2	M	299/307 (97%)	271 (91%)	27 (9%)	1 (0%)	46	84
3	H	236/260 (91%)	223 (94%)	10 (4%)	3 (1%)	15	60
All	All	813/848 (96%)	740 (91%)	64 (8%)	9 (1%)	17	63

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	H	245	ALA
1	L	201	GLU
1	L	237	SER
3	H	174	GLN
3	H	211	ASP
1	L	71	LEU
1	L	19	GLY
2	M	126	VAL
1	L	31	VAL

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.



The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	220/220 (100%)	208 (94%)	12 (6%)	27	67
2	M	236/240 (98%)	229 (97%)	7 (3%)	48	81
3	H	193/208 (93%)	188 (97%)	5 (3%)	54	83
All	All	649/668 (97%)	625 (96%)	24 (4%)	41	76

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	21	LEU
1	L	38	THR
1	L	46	ILE
1	L	154	LEU
1	L	167	PHE
1	L	185	LEU
1	L	207	ARG
1	L	216	PHE
1	L	235	LEU
1	L	247	CYS
1	L	271	TRP
1	L	272	TRP
2	M	50	ILE
2	M	148	TRP
2	M	196	LEU
2	M	204	LEU
2	M	216	PHE
2	M	223	ILE
2	M	258	PHE
3	H	105	MET
3	H	121	PRO
3	H	225	VAL
3	H	231	ASP
3	H	246	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	173	HIS
2	M	4	GLN

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Mol	Chain	Res	Type
2	M	188	ASN
2	M	259	ASN
2	M	299	GLN
3	H	129	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
5	BCL	L	851	-	53,74,74	1.35	8 (15%)	57,115,115	1.64	8 (14%)
5	BCL	L	853	-	53,74,74	1.51	7 (13%)	57,115,115	1.97	14 (24%)
6	BPH	L	855	-	64,70,70	1.25	10 (15%)	73,101,101	1.84	15 (20%)
5	BCL	M	850	-	53,74,74	1.75	7 (13%)	57,115,115	5.03	14 (24%)
5	BCL	M	852	-	53,74,74	1.39	10 (18%)	57,115,115	1.76	9 (15%)
6	BPH	M	854	-	64,70,70	1.36	10 (15%)	73,101,101	1.70	14 (19%)
7	U10	M	857	-	48,48,63	2.26	14 (29%)	58,61,79	2.34	22 (37%)
8	SPO	M	859	-	40,41,41	3.50	23 (57%)	45,50,50	2.72	15 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BCL	L	851	-	-	0/37/137/137	0/0/9/9
5	BCL	L	853	-	-	0/37/137/137	0/0/9/9
6	BPH	L	855	-	-	0/54/105/105	0/1/6/6
5	BCL	M	850	-	-	0/37/137/137	0/0/9/9
5	BCL	M	852	-	-	0/37/137/137	0/0/9/9
6	BPH	M	854	-	-	0/54/105/105	0/1/6/6
7	U10	M	857	-	-	0/45/69/87	0/1/1/1
8	SPO	M	859	-	-	0/47/47/47	0/0/0/0

All (89) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	M	857	U10	C7-C8	-3.69	1.45	1.50
8	M	859	SPO	C4-C5	-3.57	1.45	1.50
8	M	859	SPO	C31-C32	-3.29	1.41	1.50
8	M	859	SPO	C11-C12	-3.21	1.38	1.45
6	M	854	BPH	C3D-CAD	-2.79	1.41	1.46
6	L	855	BPH	C3D-CAD	-2.67	1.41	1.46
5	L	853	BCL	C3C-C4C	-2.59	1.48	1.51
6	L	855	BPH	C1B-C2B	-2.56	1.40	1.45
8	M	859	SPO	C25-C23	-2.49	1.40	1.45
6	M	854	BPH	C1B-C2B	-2.37	1.40	1.45
6	L	855	BPH	C3A-C2A	-2.31	1.47	1.54
5	M	852	BCL	CHD-C4C	-2.27	1.34	1.41
5	M	852	BCL	C3D-CAD	-2.22	1.39	1.45
5	M	850	BCL	O2D-CGD	-2.21	1.27	1.33
5	M	852	BCL	C3C-C4C	-2.19	1.48	1.51
6	L	855	BPH	O2A-CGA	-2.12	1.26	1.33
6	M	854	BPH	O2D-CGD	-2.07	1.27	1.33
5	L	851	BCL	C3C-C4C	-2.05	1.49	1.51
6	M	854	BPH	C3A-C2A	-2.04	1.48	1.54
7	M	857	U10	C12-C13	-2.02	1.44	1.50
6	L	855	BPH	C3B-CAB	2.03	1.53	1.46
8	M	859	SPO	C24-C23	2.05	1.55	1.50
5	L	853	BCL	CBB-CAB	2.07	1.55	1.49
5	M	852	BCL	CBB-CAB	2.16	1.56	1.49
6	M	854	BPH	CHA-C1A	2.18	1.42	1.37
5	M	850	BCL	OBD-CAD	2.19	1.25	1.22
5	L	851	BCL	C3B-CAB	2.23	1.55	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	M	852	BCL	CMD-C2D	2.25	1.56	1.51
6	M	854	BPH	C3B-C2B	2.26	1.44	1.38
5	L	851	BCL	CBB-CAB	2.28	1.56	1.49
6	L	855	BPH	CMD-C2D	2.34	1.56	1.51
7	M	857	U10	C8-C9	2.38	1.37	1.33
8	M	859	SPO	C22-C23	2.41	1.38	1.35
6	M	854	BPH	CMD-C2D	2.44	1.56	1.51
8	M	859	SPO	C29-C28	2.45	1.56	1.50
5	L	851	BCL	CAC-C3C	2.46	1.59	1.54
5	M	852	BCL	CMB-C2B	2.48	1.56	1.51
5	M	850	BCL	CMB-C2B	2.56	1.56	1.51
6	L	855	BPH	C3B-C2B	2.56	1.45	1.38
7	M	857	U10	C38-C39	2.60	1.40	1.32
5	L	851	BCL	CMB-C2B	2.60	1.57	1.51
5	M	852	BCL	C3B-CAB	2.76	1.56	1.49
6	L	855	BPH	C2-C3	2.78	1.38	1.33
6	M	854	BPH	C3D-C2D	2.80	1.46	1.40
8	M	859	SPO	C10-C9	2.85	1.52	1.43
6	L	855	BPH	CMB-C2B	2.86	1.56	1.50
8	M	859	SPO	C8-C7	2.87	1.57	1.50
5	L	853	BCL	CMB-C2B	2.89	1.57	1.51
8	M	859	SPO	C19-C17	2.98	1.39	1.35
7	M	857	U10	O4-C4	3.00	1.44	1.37
5	M	852	BCL	C3B-C2B	3.02	1.47	1.40
6	M	854	BPH	CMB-C2B	3.12	1.57	1.50
7	M	857	U10	C13-C14	3.14	1.39	1.33
6	L	855	BPH	C3D-C2D	3.16	1.47	1.40
5	L	853	BCL	C3B-C2B	3.21	1.47	1.40
5	L	851	BCL	C3B-C2B	3.22	1.47	1.40
8	M	859	SPO	C35-C33	3.22	1.58	1.51
5	L	853	BCL	C3B-CAB	3.27	1.57	1.49
8	M	859	SPO	C37-C38	3.28	1.42	1.32
5	L	853	BCL	C3D-C2D	3.36	1.48	1.40
5	L	851	BCL	C2-C3	3.39	1.39	1.33
7	M	857	U10	C33-C34	3.54	1.39	1.33
7	M	857	U10	C18-C19	3.63	1.40	1.33
7	M	857	U10	C6-C5	3.69	1.57	1.46
5	L	851	BCL	C3D-C2D	3.70	1.49	1.40
5	M	852	BCL	C3D-C2D	3.77	1.49	1.40
7	M	857	U10	C23-C24	3.80	1.40	1.33
5	M	852	BCL	C2-C3	3.90	1.40	1.33
5	M	850	BCL	C3D-C2D	3.91	1.49	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	M	857	U10	C28-C29	3.92	1.40	1.33
8	M	859	SPO	C15-C14	3.93	1.56	1.43
6	M	854	BPH	C2-C3	4.08	1.41	1.33
8	M	859	SPO	C13-C12	4.09	1.59	1.50
7	M	857	U10	C7-C6	4.12	1.58	1.51
7	M	857	U10	C4-C3	4.14	1.53	1.35
8	M	859	SPO	O1-CM1	4.34	1.57	1.43
8	M	859	SPO	C32-C33	4.40	1.41	1.33
8	M	859	SPO	C14-C12	4.59	1.41	1.35
5	M	850	BCL	C3B-CAB	4.67	1.61	1.49
5	M	850	BCL	C2-C3	4.79	1.42	1.33
8	M	859	SPO	C26-C25	5.02	1.47	1.34
5	L	853	BCL	C2-C3	5.46	1.43	1.33
8	M	859	SPO	C21-C20	5.73	1.51	1.35
8	M	859	SPO	C27-C28	5.78	1.40	1.34
8	M	859	SPO	C6-C5	6.86	1.51	1.31
5	M	850	BCL	C3B-C2B	6.87	1.56	1.40
8	M	859	SPO	C10-C11	6.94	1.52	1.34
7	M	857	U10	C6-C1	7.42	1.52	1.35
8	M	859	SPO	C15-C16	9.07	1.58	1.34

All (111) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	M	850	BCL	OBB-CAB-CBB	-16.77	79.93	120.13
5	M	850	BCL	CBB-CAB-C3B	-11.85	85.16	120.33
8	M	859	SPO	C25-C23-C22	-10.62	101.87	118.98
5	M	850	BCL	CMB-C2B-C1B	-8.37	114.51	128.36
5	M	852	BCL	CMB-C2B-C1B	-5.99	118.45	128.36
5	L	851	BCL	CMB-C2B-C1B	-5.93	118.55	128.36
8	M	859	SPO	C18-C17-C19	-5.87	114.23	122.90
7	M	857	U10	C15-C14-C13	-5.78	112.15	123.50
5	L	853	BCL	CMB-C2B-C1B	-5.75	118.85	128.36
5	L	853	BCL	CAA-C2A-C1A	-5.54	92.94	112.47
7	M	857	U10	C10-C9-C8	-5.21	113.27	123.50
8	M	859	SPO	C24-C23-C22	-5.09	115.38	122.90
8	M	859	SPO	C20-C21-C22	-4.84	112.69	123.39
6	M	854	BPH	O1D-CGD-CBD	-4.79	117.75	124.62
5	M	850	BCL	CAA-C2A-C1A	-4.61	96.20	112.47
6	L	855	BPH	O1D-CGD-CBD	-4.45	118.25	124.62
7	M	857	U10	C35-C34-C33	-4.31	115.03	123.50
6	L	855	BPH	C4D-C3D-C2D	-4.15	101.72	107.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	M	854	BPH	C4D-C3D-C2D	-4.06	101.84	107.08
6	L	855	BPH	C5-C3-C2	-3.89	113.67	121.05
7	M	857	U10	C25-C24-C23	-3.81	116.02	123.50
5	M	852	BCL	OBD-CAD-C3D	-3.80	120.60	128.35
8	M	859	SPO	C15-C14-C12	-3.78	121.73	127.20
8	M	859	SPO	C4-C5-C6	-3.71	119.38	124.67
6	L	855	BPH	CHC-C4B-NB	-3.70	117.87	124.91
7	M	857	U10	C20-C19-C18	-3.66	116.32	123.50
5	L	853	BCL	OBD-CAD-C3D	-3.65	120.90	128.35
6	L	855	BPH	CMB-C2B-C1B	-3.54	119.30	125.06
6	M	854	BPH	CHC-C4B-NB	-3.48	118.28	124.91
8	M	859	SPO	C11-C12-C14	-3.41	113.49	118.98
5	M	850	BCL	O1D-CGD-CBD	-3.29	119.90	124.62
5	M	850	BCL	OBD-CAD-C3D	-3.20	121.82	128.35
5	L	851	BCL	OBD-CAD-C3D	-3.11	122.01	128.35
6	M	854	BPH	CMB-C2B-C1B	-3.09	120.03	125.06
8	M	859	SPO	C15-C16-C17	-2.91	117.76	126.32
5	L	853	BCL	O1D-CGD-CBD	-2.90	120.47	124.62
7	M	857	U10	C31-C32-C33	-2.81	104.32	111.69
6	M	854	BPH	CBB-CAB-C3B	-2.79	114.33	120.52
5	M	852	BCL	CHA-C1A-NA	-2.72	119.37	126.06
5	L	853	BCL	CAC-C3C-C2C	-2.70	107.34	114.13
7	M	857	U10	O5-C5-C4	-2.62	115.11	120.79
5	L	851	BCL	CHA-C1A-NA	-2.44	120.05	126.06
5	L	853	BCL	C16-C15-C13	-2.34	107.72	115.49
7	M	857	U10	C30-C29-C28	-2.33	118.93	123.50
7	M	857	U10	C1-C6-C5	-2.30	117.50	120.12
8	M	859	SPO	C9-C10-C11	-2.29	116.16	123.13
6	L	855	BPH	C1C-NC-C4C	-2.19	108.19	110.44
8	M	859	SPO	C34-C33-C35	-2.13	112.15	115.41
5	M	850	BCL	CHA-C1A-NA	-2.13	120.82	126.06
6	L	855	BPH	CBB-CAB-C3B	-2.13	115.80	120.52
6	M	854	BPH	C4-C3-C2	-2.11	119.35	123.50
6	M	854	BPH	C1C-NC-C4C	-2.09	108.30	110.44
5	M	852	BCL	C15-C13-C12	-2.05	100.02	112.27
7	M	857	U10	C27-C28-C29	-2.03	123.36	127.76
8	M	859	SPO	C10-C9-C7	-2.01	124.29	127.20
5	L	851	BCL	C3D-CAD-CBD	2.03	110.46	107.60
5	L	853	BCL	C2C-C3C-C4C	2.03	104.94	101.50
5	L	853	BCL	C3D-CAD-CBD	2.08	110.53	107.60
6	L	855	BPH	C2A-C3A-C4A	2.20	106.13	101.10
7	M	857	U10	C10-C9-C11	2.21	118.78	115.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	L	855	BPH	CED-O2D-CGD	2.26	121.29	115.99
6	M	854	BPH	CMB-C2B-C3B	2.31	133.41	128.04
6	M	854	BPH	CED-O2D-CGD	2.32	121.43	115.99
6	L	855	BPH	C3D-CAD-CBD	2.39	110.97	107.60
5	L	851	BCL	C2A-C1A-CHA	2.40	128.30	123.89
5	L	853	BCL	O2D-CGD-CBD	2.41	114.60	111.30
5	M	850	BCL	O2A-CGA-CBA	2.43	119.30	111.90
5	L	853	BCL	O2A-CGA-CBA	2.44	119.33	111.90
7	M	857	U10	C35-C34-C36	2.45	119.15	115.41
6	L	855	BPH	CMB-C2B-C3B	2.50	133.87	128.04
7	M	857	U10	C36-C34-C33	2.51	125.81	121.05
5	M	850	BCL	C2A-C1A-CHA	2.52	128.53	123.89
5	M	852	BCL	C3D-CAD-CBD	2.54	111.19	107.60
6	M	854	BPH	C6-C5-C3	2.56	118.10	112.48
8	M	859	SPO	C16-C17-C19	2.64	123.24	118.98
8	M	859	SPO	C18-C17-C16	2.66	122.53	118.10
5	M	852	BCL	C2A-C1A-CHA	2.66	128.79	123.89
8	M	859	SPO	O1-C1-C4	2.67	112.40	105.87
5	L	851	BCL	C6-C5-C3	2.72	118.44	112.48
7	M	857	U10	C11-C12-C13	2.83	119.10	111.69
5	L	853	BCL	C6-C5-C3	2.85	118.73	112.48
6	M	854	BPH	C3D-CAD-CBD	2.86	111.64	107.60
7	M	857	U10	C30-C29-C31	2.87	119.79	115.41
7	M	857	U10	C21-C19-C18	3.01	126.76	121.05
7	M	857	U10	C4M-O4-C4	3.03	127.39	116.61
5	M	850	BCL	CBA-CAA-C2A	3.19	122.74	113.73
7	M	857	U10	C25-C24-C26	3.23	120.34	115.41
5	M	850	BCL	O2D-CGD-CBD	3.26	115.77	111.30
8	M	859	SPO	C8-C7-C9	3.45	128.00	122.90
5	M	850	BCL	C6-C5-C3	3.57	120.31	112.48
7	M	857	U10	C11-C9-C8	3.63	127.94	121.05
7	M	857	U10	C7-C8-C9	3.64	132.86	126.70
5	M	850	BCL	CED-O2D-CGD	3.78	124.86	115.99
6	M	854	BPH	C4-C3-C5	3.87	121.32	115.41
5	L	853	BCL	CMB-C2B-C3B	4.07	133.05	125.09
5	L	853	BCL	CBA-CAA-C2A	4.13	125.40	113.73
5	L	851	BCL	CED-O2D-CGD	4.22	125.89	115.99
5	M	852	BCL	CMB-C2B-C3B	4.36	133.61	125.09
5	M	852	BCL	CED-O2D-CGD	4.38	126.26	115.99
5	L	853	BCL	CED-O2D-CGD	4.45	126.42	115.99
5	L	851	BCL	CMB-C2B-C3B	4.47	133.82	125.09
6	L	855	BPH	O2D-CGD-CBD	4.54	117.52	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	L	855	BPH	C3C-C4C-NC	4.78	112.72	107.93
6	M	854	BPH	O2D-CGD-CBD	4.80	117.89	111.30
6	L	855	BPH	C6-C5-C3	4.85	123.13	112.48
5	M	852	BCL	C6-C5-C3	4.91	123.27	112.48
6	M	854	BPH	C3C-C4C-NC	4.93	112.87	107.93
7	M	857	U10	C15-C14-C16	5.48	123.78	115.41
7	M	857	U10	C7-C6-C5	5.56	125.09	118.56
6	L	855	BPH	C4-C3-C5	5.89	124.40	115.41
5	M	850	BCL	OBB-CAB-C3B	28.46	165.09	120.00

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 44 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	851	BCL	11	0
5	L	853	BCL	5	0
6	L	855	BPH	8	0
5	M	850	BCL	11	0
5	M	852	BCL	6	0
6	M	854	BPH	9	0
7	M	857	U10	3	0
8	M	859	SPO	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.